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Fax

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Fax:	703 746 5318		Date:	12 June 2003	
Phone:			Pages:	(including cover	(including cover sheet) 12
Your Ref.:	App #09/846,259	1	Our Ref.:	0425-0819P	
Re:	Our 10 June 2003 Conference		CC:		
Urgent	☑ For Review	☑ Please Comment		Please Reply	Please Recycle

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Comments:

Thank you very much for telephoning our offices on 10 June 2003. It is our understanding, based on that conference, that if the application were amended as shown on the attached sheets that all outstanding issues would be resolved and that you would pass the case to issue. No prior art was discussed.

If our understanding is correct please so inform us by any means convenient to you. We reliably receive e-mail sent to mailroom@bskb.com.

As soon as we have your response we will contact our client, who will determine whether, and in what manner the application will be amended.

Dand M. Muyhy

App \$ 09/846,259 Dk+ 0425-0189P

What is claimed is:

(currently amendal)

1. A 1,4-substituted cyclic amine derivative represented by the compound of following formula (I):

$$R^{1}$$
 $B-C$ R^{2} D $CCH_{2})_{m}$ $T-(CH_{2})_{m}$ Y $Z-R$ $CCH_{2})_{p}$ $CCH_{2})_{p}$

wherein A, B, C, D, and T are the same or different from one another and each represents methine or nitrogen, provided that one and only one of them represents nitrogen;

the bond represented by the following formula:

represents a single or double bond;

Y and Z are the same or different from each other and each represents methine, nitrogen, a group represented by the following formula:

or a group represented by the following formula:

provided at least one of them represents nitrogen;

R¹ and R² are the same or different from each other and each represents

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hydrogen, halogeno, hydroxy, lower alkylsulfonylaminoalkyl, lower halogenatedalkylsulfonylaminoalkyl, 2-pyrrolidinon-1-yl, 1-hydroxy-1-(methoxypyridyl)methyl, methoxypyridylcarbonyl, 1,3-propanesultum-2-yl, lower hydroxypiperidylcarbonylalkyl, lower hydroxyalkylamidoalkyl, lower halogenated-alkylamidoalkyl, lower dihalogenatedalkylamidoalkyl, lower heteroarylamidoalkyl, lower hydroxyalkylamidoalkyl, optionally substituted amino, nitro, lower alkyl, lower alkoxy, lower acyl, lower alkoxyalkoxy, cyano, lower alkylsulfonyl, sulfonylamido, hydroxy-lower alkyl, hydroxy-lower alkoxy, lower alkoxycarbonylamino, lower alkylsulfonylamino, N-lower alkylalkylsulfonylamino, lower acylamino, optionally substituted aminoalkyl, optionally N-substituted lower acylaminoalkyl, optionally substituted aryl, optionally substituted arylsulfonylamino, lower alkylsulfonyloxy, hydroxyiminomethyl, (2-pyrrolidon1-yl)methyl, (2-piperidon-1-yl)methyl, optionally substituted heteroaryl, optionally substituted aralkyl, optionally substituted heteroarylalkyl, cycloalkylcarbonylaminoalkyl, optionally substituted ureido, optionally substituted ureido-lower alkyl, succinimido, (succinimido-1-yl)lower alkyl, amido, optionally substituted carbamoyl, optionally substituted carbamoyl-lower alkyl, optionally substituted thiocarbamoyllower alkyl, formyl, aromatic acyl, heteroarylcarbonyl, halogenated lower alkyl, (2-imidazolidinon -1-yl)methyl, (2,4-imidazolidinedion-3-yl)methyl, (2-oxazolidon3-yl)methyl, (glutarimido-1-yl)methyl, optionally substituted heteroarylhydroxyalkyl, cyano-lower alkyl, 1-hydroxy lower cycloalkyl, (2,4-thiazolidinedion-3-yl)methyl, optionally substituted 4-piperidylmethyl, heteroarylacyl, pyrrolidinylcarbonyl-lower alkyl, optionally substituted aminosulfonylalkyl, carboxy-lower alkyl, or lower alkylamidoalkyl; or alternatively R¹ and R² together may form optionally substituted alicycle) optionally substituted heterocycle or alkylenedioxy, provided

0/m

these rings may be substituted;

R³ represents hydrogen, halogeno, lower alkyl, hydroxy, hydroxy-lower alkyl, lower alkoxy, formyl, optionally substituted aralkyloxy, hydroxy-lower alkoxy, optionally substituted sulfamoyl, or optionally N-substituted sulfamoyl-lower alkyl;

R⁴ represents hydrogen, lower alkyl, hydroxy-lower alkyl, lower alkoxyalkyl, optionally aryl-substituted aryloxyalkyl, or optionally aryl-substituted aralkyloxyalkyl;

R⁵ represents lower alkyl, lower acyl, lower alkoxycarbonyl, aromatic acyl, or a group represented by the following formula:

$$-Q^{1}-(CH_{2})_{S}-Q^{2}-R^{6}$$

wherein Q^1 and Q^2 are both single bonds, or one of them is a single bond while the other represents oxygen, carbonyl, a group represented by -NHCO-, a group represented by -NHSO₂-, or a group represented by >CH-R⁷, wherein R⁷ represents hydroxy, lower alkyl or halogeno:

s represents 0 or an integer of 1 to 6; and

R⁶ represents optionally substituted aryl, optionally substituted heteroaryl, optionally substituted benzoheteroaryl) 1,4-benzodioxanyl, 1,3-benzodioxolyl, benzothiazolyl, or cyano;

n represents 1;

m represents 0 or an integer of 1 to 6; and

p represents an integer of 1 to 3,

and pharmacologically acceptable salts thereof.

8/m

(currently emended)

2. A 1-74-substituted cyclic amir compound of following formula:

wherein R represents a substituent of the formula:

$$\begin{pmatrix} R^1 & R^2 \\ R^3 & R^2 \end{pmatrix}$$
 or
$$\begin{pmatrix} R^1 & R^2 \\ R^3 & R^2 \end{pmatrix}$$

wherein the bond represented by the following formula:

and R^1 , R^2 , R^3 , R^4 , R^5 , Y, Z, m, and p are each as defined in claim 1, and pharmacologically acceptable salts thereof.

(currently amended) compound of

3. The +4-substituted cyclic amine derivative as set forth in claim 1 or a pharmacologically acceptable salt thereof, wherein m is 0 and p is 2.

a pharmacologically acceptable salt thereof, wherein Y is methine and Z is nitrogen.

(currently amended) compound of The 1,4 substituted cyclic amine derivative as set forth in in claim 1

5. or a pharmacologically acceptable salt thereof, which is a compound selected from among the following ones:

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- 1-{1-[2-(4-methoxyphenyl)ethyl]piperidin-4-yl}-7-methoxy-1,2,3,4-tetrahydroquinoline,
- (268) 1-{1-[2-(4-fluorophenyl)ethyl]piperidin-4-yl}-7-methoxy-1,2,3,4-tetrahydroquinoline,
- (269) 1-[1-(4-cyanopropyl)piperidin-4-yl]-7-methoxy-1,2,3,4-tetrahydroquinoline, (270)
- 1-{1-[2-(2-thienyl)ethyl]piperidin-4-yl}-7-methoxy-1,2,3,4-tetrahydroquinoline, (271)
- 1-{1-[2-(4-fluorophenyl)ethyl]piperidin-4-yl}-7,8-dimethoxy-1,2,3,4-tetrahydroquinoline,
- (272) 1-{1-[2-(4-fluorophenyl)ethyl]piperidin-4-yl}-7,8-methylenedioxy-1,2,3,4tetrahydroguinoline,
- (273) 1-{1-[2-(4-fluorophenyl)ethyl]piperidin-4-yl}-7-methoxy-8-methyl-1,2,3,4tetrahydroquinoline,
- (274) 1-{1-[2-(4-fluorophenyl)-2-oxoethyl]piperidin-4-yl}-7-methoxy-1,2,3,4tetrahydroquinoline,
- (275) 1-{1-[2-(4-fluorophenyl)-2-hydroxyethyl]piperidin-4-yl}-7-methoxy-1,2,3,4tetrahydroguinoline,
- (276) 1-{1-[2-(4-fluorophenyl)-2-fluoroethyl]piperidin-4-yl}-7-methoxy-1,2,3,4tetrahydroquinoline, and
- (283) 5-{4-[2-(4-fluorophenyl)ethyl]piperazin-1-yl}-5,6,7,8-tetrahydroisoquinoline.

(currently amended)

6. A pharmaceutical composition comprising a therapeutically effective compound of amount of the 1,4 substituted cyclic amine derivative or salt as set forth in claim 1 or its salt

7. An agent for treating ameliorating, and preventing diseases against which serotonin antagonism is efficacious, which contain as the active ingredient the 1,4 substituted cyclic amine derivative as set forth in claim 1 or a pharmacologically acceptable salt thereof.

(currently emends)

8. An agent for treating, ameliorating, and preventing spastic paralysis, which contain as the active amount of the compound of which contain as the active ingredient, the 1,1 substituted cyclic amine derivative as set forth in claim 1 or a pharmacologically acceptable salt thereof.

9. A muscle relaxant which contains as the active ingredient the an effective amount of the compound of 1,4 substituted cyclic amine derivative as set forth in claim 1 or a in a mixture pharmacologically acceptable salt thereof.

10. A process for producing a 1,4-substituted cyclic amine derivative compand of the represented by the following formula:

wherein the bond represented by the following formula:

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and A, B, C, D, R¹, R², R³, R⁴, R⁵, n, and p are each as defined in claim 1, which comprises removing, if necessary, the protecting group from a 1,4-substituted cyclic amine derivative (IX) represented by the following formula:

$$(CH_2)_n = R^2$$

$$(CH_2)_n = N$$

$$R^4$$

$$(CH_2)_n = N$$

$$R^3$$

$$(CH_2)_n = N$$

$$R^4$$

$$(CH_2)_n = N$$

wherein the bond represented by the following formula:

and A, B, C, D, R^1 , R^2 , R^3 , R^4 , n, and p are each as defined in claim 1; and Pr.G represents hydrogen or a protecting group, and then reacting the same with L-R⁵ wherein R^5 is as defined in claim 1; and L represents a leaving group.

11. A process for producing 4,4-substituted cyclic amine derivative (X), a compound of as set forth in claim 1, which comprises reacting a fused cyclic amine represented by the following formula:

wherein the bond represented by the following formula:

and A, B, C, D, R¹, R², R³ and n are each as defined in claim 1 with a cyclic ketone (VIII) represented by the following formula:

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$$O = \begin{pmatrix} R^4 \\ N-Pr.G \\ (CH_2)_p \end{pmatrix}$$
 (VIII)

wherein R4, p, and Pr.G are each as defined in claim 1

in the presence of a reducing agent to thereby give a 1,4-substituted cyclic amine derivative (IX), removing, if necessary, the protecting group therefrom and further reacting the same with L-R⁵.

12. A 4-substituted cyclic amine derivative represented by the following formula:

wherein the bond represented by the following formula:

and A, B, C, D, R^1 , R^2 , R^3 , R^4 , n, and p are each as defined in claim 1, provided that the case where R^1 , R^2 , R^3 and R^4 are all hydrogen atoms is excluded.

(currently amendal)

(currently amendal)

13. A method for treating a disease to which serotonin antagonism is

efficacious, which comprises administering an effective dose of the 1,4-substituted cyclic amine derivative as set forth-in claim 1, or a pharmacologically acceptable salt thereof, to a person in need of such treatment.

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(currently amendal)

14. The 1,4-substituted cyclic amine derivative as set forth in claim 1, in which the bond represented by the following formula in the formula (I):

is a single bond, represented by the formula (XXI):

or a pharmacologically acceptable salt thereof.

(currently amended)

15. The 4,4 substituted cyclic amine derivative as set forth-in claim 1, in which m is 0 in the formula (I), represented by the formula (XXII):

or a pharmacologically acceptable salt thereof.

16. The 1,4 substituted cyclic amine derivative as set forth in Claim 1, in which m is 1 to 6 in the formula (I) or a pharmacologically acceptable salt thereof.

(12/12/2) (12/12/2)

compound of decivative represented by the formula

:(IIIXX)

or a pharmacologically acceptable salt thereof.

(arrestly amount of the 18. The 1.4 substituted cyclic amine derivative as set forth in claim 1, in

which the bond represented by the following formula in the formula (I):

is a double bond, represented by the formula (XXIV):

or a pharmacologically acceptable salt thereof.

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19. The 1,4 substituted cyclic amine derivative as set forth in claim 1, in which the T is nitrogen.

ABSTRACT OF THE DISCLOSURE

A novel 1,4-substituted cyclic amine derivative represented by the compound of the following formula:

$$R^{1}$$
 $B-C$ R^{2}
 $(CH_{2})_{n}$ $T-(CH_{2})_{m}-Y$ $Z-R^{5}$
 $(CH_{2})_{0}$

wherein A, B, C, D, T, Y, and Z represent each methine or nitrogen; R¹, R², R³, R⁴, and R⁵ represent each a substituent; n represents 0 or an integer of 1 to 3; m represents 0 or an integer of 1 to 6; and p represents an integer of 1 to 3; and pharmacologically acceptable salts thereof. The compound has a serotonin antagonism and is clinically useful as medicament, in particular, for treating, ameliorating and preventing spastic paralysis or central muscle relaxants for ameliorating myotonia.

